

| | Final Report |
|----------|----------------|
| √ | Revised Report |

Report Date: 14-Jan-19 12:59

Laboratory Report SC52773

Gulf Oil L.P. 281 Eastern Avenue Chelsea, MA 02150 Attn: Andrew P. Adams

Project: Gulf Terminal - Chelsea, MA

Project #: Gulf Chelsea

I attest that the information contained within the report has been reviewed for accuracy and checked against the quality control requirements for each method. These results relate only to the sample(s) as received.

All applicable NELAC requirements have been met.

Massachusetts # M-MA138/MA1110 Connecticut # PH-0777 Florida # E87936 Maine # MA138 New Hampshire # 2972/2538 New Jersey # MA011 New York # 11393 Pennsylvania # 68-04426/68-02924 Rhode Island # LAO00348 USDA # P330-15-00375 Vermont # VT-11393



Authorized by:

Dawn Wojcik Laboratory Director



Eurofins Spectrum Analytical holds primary certification in the State of Massachusetts for the analytes as indicated with an X in the "Cert." column within this report. Please note that the State of Massachusetts does not offer certification for all analytes. Please refer to our website for specific certification holdings in each state.

Please note that this report contains 24 pages of analytical data plus Chain of Custody document(s). When the Laboratory Report is indicated as revised, this report supersedes any previously dated reports for the laboratory ID(s) referenced above. Where this report identifies subcontracted analyses, copies of the subcontractor's test report are available upon request. This report may not be reproduced, except in full, without written approval from Eurofins Spectrum Analytical, Inc.

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Please contact the Laboratory or Technical Director at 800-789-9115 with any questions regarding the data contained in this laboratory report.

Sample Summary

Work Order: SC52773

Project: Gulf Terminal - Chelsea, MA

Project Number: Gulf Chelsea

| Laboratory ID | Client Sample ID | <u>Matrix</u> | Date Sampled | Date Received |
|---------------|------------------|---------------|-----------------|----------------------|
| SC52773-01 | Outfall 003 | Surface Water | 19-Dec-18 13:15 | 20-Dec-18 16:30 |
| SC52773-02 | Trip Blank | Trip Blank | 19-Dec-18 00:00 | 20-Dec-18 16:30 |

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CASE NARRATIVE:

Data has been reported to the RDL. This report excludes estimated concentrations detected below the RDL and above the MDL (J-Flag).

All non-detects and all results below the reporting limit are reported as "<" (less than) the reporting limit in this report.

The samples were received 1.6 degrees Celsius, please refer to the Chain of Custody for details specific to temperature upon receipt. An infrared thermometer with a tolerance of +/- 1.0 degrees Celsius was used immediately upon receipt of the samples.

If a Matrix Spike (MS), Matrix Spike Duplicate (MSD) or Duplicate (DUP) was not requested on the Chain of Custody, method criteria may have been fulfilled with a source sample not of this Sample Delivery Group. If method or program required MS/MSD/Dup were not performed, sufficient sample was not provided to the laboratory.

Analyses for Total Hardness, pH, and Total Residual Chlorine fall under the state of Pennsylvania code Chapter 252.6 accreditation by rule.

January 14, 2019 Report Revision Case Narrative:

This report has been revised to remove 2-Methylnaphthalene.

See below for any non-conformances and issues relating to quality control samples and/or sample analysis/matrix.

E1664A

Laboratory Control Samples:

CC19445-LCS

A Blank spike was performed instead of a matrix spike

Oil and Grease by EPA 1664A

CC19445-BLK

A Blank spike was performed instead of a matrix spike

Oil and Grease by EPA 1664A

CC19445-LCSD

A Blank spike was performed instead of a matrix spike

Oil and Grease by EPA 1664A

SW8260C

Laboratory Control Samples:

CC19490-LCS

Laboratory Control Samples:

CC19490-LCS

A LCS and LCS Duplicate were performed instead of a matrix spike and matrix spike duplicate.

- % 1,2-dichlorobenzene-d4
- % Bromofluorobenzene
- % Dibromofluoromethane
- % Toluene-d8
- 1,1,1,2-Tetrachloroethane
- 1,1,1-Trichloroethane
- 1,1,2,2-Tetrachloroethane
- 1,1,2-Trichloroethane
- 1,1-Dichloroethane
- 1,1-Dichloroethene
- 1,1-Dichloropropene
- 1,2,3-Trichlorobenzene
- 1,2,3-Trichloropropane
- 1,2,4-Trichlorobenzene
- 1,2,4-Trimethylbenzene
- 1,2-Dibromo-3-chloropropane
- 1,2-Dibromoethane
- 1,2-Dichlorobenzene
- 1,2-Dichloroethane
- 1,2-Dichloropropane
- 1,3,5-Trimethylbenzene
- 1,3-Dichlorobenzene
- 1,3-Dichloropropane 1,4-Dichlorobenzene
- 1,4-dioxane
- 2,2-Dichloropropane
- 2-Chlorotoluene
- 2-Hexanone
- 2-Isopropyltoluene
- 4-Chlorotoluene
- 4-Methyl-2-pentanone

Acetone

Acrylonitrile

Benzene

Bromobenzene

Bromochloromethane

Bromodichloromethane

Bromoform

Bromomethane

Carbon Disulfide

Carbon tetrachloride

Chlorobenzene Chloroethane

Chloroform

Chloromethane

cis-1,2-Dichloroethene

cis-1,3-Dichloropropene

Dibromochloromethane

Dibromomethane

Dichlorodifluoromethane

Ethyl ether

Ethylbenzene

Hexachlorobutadiene

Isopropylbenzene

m&p-Xylene

Methyl ethyl ketone

Laboratory Control Samples:

CC19490-LCS

A LCS and LCS Duplicate were performed instead of a matrix spike and matrix spike duplicate.

Methyl t-butyl ether (MTBE)

Methylene chloride

Naphthalene

n-Butylbenzene

n-Propylbenzene

o-Xylene

p-Isopropyltoluene

sec-Butylbenzene

Styrene

tert-Butylbenzene

Tetrachloroethene

Tetrahydrofuran (THF)

Toluene

trans-1,2-Dichloroethene

trans-1,3-Dichloropropene

trans-1,4-dichloro-2-butene

Trichloroethene

Trichlorofluoromethane

Trichlorotrifluoroethane

Vinyl chloride

CC19490-BLK

CC19490-BLK

A LCS and LCS Duplicate were performed instead of a matrix spike and matrix spike duplicate.

- % 1,2-dichlorobenzene-d4
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- 1,1,1-Trichloroethane
- 1,1,2,2-Tetrachloroethane
- 1,1,2-Trichloroethane
- 1,1-Dichloroethane
- 1,1-Dichloroethene
- 1,1-Dichloropropene
- 1,2,3-Trichlorobenzene
- 1,2,3-Trichloropropane
- 1,2,4-Trichlorobenzene
- 1,2,4-Trimethylbenzene
- 1,2-Dibromo-3-chloropropane
- 1,2-Dibromoethane
- 1,2-Dichlorobenzene
- 1,2-Dichloroethane
- 1,2-Dichloropropane
- 1,3,5-Trimethylbenzene
- 1,3-Dichlorobenzene
- 1,3-Dichloropropane
- 1,4-Dichlorobenzene
- 1,4-dioxane
- 2,2-Dichloropropane
- 2-Chlorotoluene
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- 2-Isopropyltoluene
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Acetone

Acrylonitrile

Benzene

Bromobenzene

Bromochloromethane

Bromodichloromethane

Bromoform

Bromomethane

Carbon Disulfide

Carbon tetrachloride

Chlorobenzene

Chloroethane

Chloroform

Chloromethane

cis-1,2-Dichloroethene

cis-1,3-Dichloropropene

Dibromochloromethane

Dibromomethane

Dichlorodifluoromethane

Ethyl ether

Ethylbenzene

Hexachlorobutadiene

Isopropylbenzene

m&p-Xylene

Methyl ethyl ketone

Methyl t-butyl ether (MTBE)

Methylene chloride

CC19490-BLK

A LCS and LCS Duplicate were performed instead of a matrix spike and matrix spike duplicate.

Naphthalene

n-Butylbenzene

n-Propylbenzene

o-Xylene

p-Isopropyltoluene

sec-Butylbenzene

Styrene

tert-Butylbenzene

Tetrachloroethene

Tetrahydrofuran (THF)

Toluene

trans-1,2-Dichloroethene

trans-1,3-Dichloropropene

trans-1,4-dichloro-2-butene

Trichloroethene

Trichlorofluoromethane

Trichlorotrifluoroethane

Vinyl chloride

CC19490-LCSD

CC19490-LCSD

A LCS and LCS Duplicate were performed instead of a matrix spike and matrix spike duplicate.

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- % Bromofluorobenzene
- % Dibromofluoromethane
- % Toluene-d8
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- 1,1,1-Trichloroethane
- 1,1,2,2-Tetrachloroethane
- 1,1,2-Trichloroethane
- 1,1-Dichloroethane
- 1,1-Dichloroethene
- 1,1-Dichloropropene
- 1,2,3-Trichlorobenzene
- 1,2,3-Trichloropropane
- 1,2,4-Trichlorobenzene
- 1,2,4-Trimethylbenzene
- 1,2-Dibromo-3-chloropropane
- 1,2-Dibromoethane
- 1,2-Dichlorobenzene
- 1,2-Dichloroethane
- 1,2-Dichloropropane
- 1,3,5-Trimethylbenzene
- 1,3-Dichlorobenzene
- 1,3-Dichloropropane 1,4-Dichlorobenzene
- 1,4-dioxane
- 2,2-Dichloropropane
- 2-Chlorotoluene
- 2-Hexanone
- 2-Isopropyltoluene
- 4-Chlorotoluene
- 4-Methyl-2-pentanone

Acetone

Acrylonitrile

Benzene

Bromobenzene

Bromochloromethane

Bromodichloromethane

Bromoform

Bromomethane

Carbon Disulfide

Carbon tetrachloride

Chlorobenzene

Chloroform

Cinorororini

Chloromethane

cis-1,2-Dichloroethene

cis-1,3-Dichloropropene

Dibromochloromethane

Dibromomethane

Dichlorodifluoromethane

Ethyl ether

Ethylbenzene

Hexachlorobutadiene

Isopropylbenzene

m&p-Xylene

Methyl ethyl ketone

Methyl t-butyl ether (MTBE)

Methylene chloride

CC19490-LCSD

A LCS and LCS Duplicate were performed instead of a matrix spike and matrix spike duplicate.

Naphthalene

n-Butylbenzene

n-Propylbenzene

o-Xylene

p-Isopropyltoluene

sec-Butylbenzene

Styrene

tert-Butylbenzene

Tetrachloroethene

Tetrahydrofuran (THF)

Toluene

trans-1,2-Dichloroethene

trans-1,3-Dichloropropene

trans-1,4-dichloro-2-butene

Trichloroethene

Trichlorofluoromethane

Trichlorotrifluoroethane

Vinyl chloride

SW8260C.

Laboratory Control Samples:

CC19490-LCS

A blank MS/MSD was analyzed with this batch.

Ethanol

CC19490-BLK

A blank MS/MSD was analyzed with this batch.

Ethanol

CC19490-LCSD

A blank MS/MSD was analyzed with this batch.

Ethanol

CC19490-MS

A blank MS/MSD was analyzed with this batch.

Ethanol

CC19490-MSD

A blank MS/MSD was analyzed with this batch.

Ethanol

SW8270D (SIM)

Laboratory Control Samples:

CC17546-LCS

This parameter is outside laboratory lcs/lcsd specified recovery limits.

% Nitrobenzene-d5

Naphthalene

SW8270D (SIM)

Laboratory Control Samples:

CC17546-LCS

This parameter is outside laboratory rpd specified recovery limits.

% Nitrobenzene-d5

Naphthalene

CC17546-LCSD

This parameter is outside laboratory lcs/lcsd specified recovery limits.

% Nitrobenzene-d5

2-Methylnaphthalene

Naphthalene

This parameter is outside laboratory rpd specified recovery limits.

% Nitrobenzene-d5

Naphthalene

CC17546-MS

This parameter is outside laboratory rpd specified recovery limits.

% 2-Fluorobiphenyl

% Nitrobenzene-d5

Acenaphthene

Benzo(ghi)perylene

Dibenz(a,h)anthracene

Fluoranthene

Fluorene

Indeno(1,2,3-cd)pyrene

Naphthalene

Pyrene

CC17546-MSD

This parameter is outside laboratory ms/msd specified recovery limits.

Benzo(ghi)perylene

Dibenz(a,h)anthracene

Indeno(1,2,3-cd)pyrene

This parameter is outside laboratory rpd specified recovery limits.

% 2-Fluorobiphenyl

% Nitrobenzene-d5

Acenaphthene

Benzo(ghi)perylene

Dibenz(a,h)anthracene

Fluoranthene

Fluorene

Indeno(1,2,3-cd)pyrene

Naphthalene

Pyrene

Sample Acceptance Check Form

| Project: | Gulf Terminal - Chelsea, MA / Gulf Chelsea | | | |
|---------------------------|---|--------------|-----------|--------------|
| Work Order: | SC52773 | | | |
| Sample(s) received on: | 12/20/2018 | | | |
| | | | | |
| The following outlines to | he condition of samples for the attached Chain of Custody upon receipt. | | | |
| | | <u>Yes</u> | <u>No</u> | <u>N/A</u> |
| Were custody se | als present? | | ✓ | |
| Were custody se | als intact? | | | \checkmark |
| Were samples re | seeived at a temperature of $\leq 6^{\circ}$ C? | \checkmark | | |
| Were samples re | frigerated upon transfer to laboratory representative? | ✓ | | |
| Were sample co | ntainers received intact? | \checkmark | | |
| | roperly labeled (labels affixed to sample containers and include sample ID, site project number and the collection date)? | | | |
| Were samples ac | ecompanied by a Chain of Custody document? | \checkmark | | |
| include sample l | Custody document include proper, full, and complete documentation, which shall (D, site location, and/or project number, date and time of collection, collector's name, e, sample matrix and any special remarks concerning the sample? | | | |
| Did sample cont | ainer labels agree with Chain of Custody document? | \checkmark | | |
| Were samples re | ceived within method-specific holding times? | abla | П | |

Client:

Gulf Oil L.P.

Summary of Hits

Lab ID: SC52773-01

| Parameter | Result | Flag | Reporting Limit | Units | Analytical Method |
|-----------------------------|--------|------|-----------------|-------|-------------------|
| Oil and Grease by EPA 1664A | < 1.5 | | 1.5 | mg/L | E1664A |
| Total Suspended Solids | 52 | | 5.0 | mg/L | SM2540D-11 |

Client ID:

Outfall 003

Please note that because there are no reporting limits associated with hazardous waste characterizations or micro analyses, this summary does not include hits from these analyses if included in this work order.

| Sample Id Outfall 00 SC52773- | | | | Client P Gulf C | | | Matrix Surface Wa | | ection Date -Dec-18 13 | | | ceived Dec-18 | |
|-------------------------------------|---|--------------------|--------------|--------------------|-------|------|----------------------|--------------------|------------------------|--------------------|---------|------------------|------|
| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | MDL | Dilution | Method Ref. | Prepared | Analyzed | Analyst | Batch | Cert |
| General C | hemistry Parameters | | | | | | | | | | | | |
| | рН | 7.03 | рН | pH Units | | | 1 | ASTM D 1293-99B | 21-Dec-18 11:30 | 21-Dec-18 17:58 | BD | 1816206 | X |
| | cted Analyses by method E1664A | | | | | | | | | | | | |
| Analysis pe | erformed by Phoenix Environi | nental Labs, | Inc. * - MAC | T007 | | | | | | | | | |
| | Oil and Grease by EPA 1664A | < 1.5 | | mg/L | 1.5 | 1.5 | 1.1 | E1664A | 19-Dec-18 13:15 | 27-Dec-18 07:54 | M-CT007 | 461365A | i |
| | by method SM2540D-11 | | | | | | | | | | | | |
| Analysis pe | erformed by Phoenix Environi Total Suspended Solids | nental Labs, . 52 | Inc. * - MAC | mg/L | 5.0 | 5.0 | 1 | SM2540D-11 | " | 26-Dec-18 07:23 | M-CT007 | 461248A | ı |
| Prepared | by method SW8260C | u outal I aba | Luo * MAC | T007 | | | | | | | | | |
| Anaiysis pe 71-43-2 | erformed by Phoenix Environi Benzene | < 0.70 | inc. · - MAC | ug/L | 0.70 | 0.70 | 1 | SW8260C | " | 24-Dec-18 11:44 | M-CT007 | 461279A | ı |
| 1634-04-4 | Methyl t-butyl ether (MTBE) | < 1.0 | | ug/L | 1.0 | 1.0 | 1 | " | " | " | " | " | |
| 91-20-3 | Naphthalene | < 1.0 | | ug/L | 1.0 | 1.0 | 1 | H . | " | " | " | " | |
| Surrogate i | recoveries: | | | | | | | | | | | | |
| 2199-69-1 | % 1,2-dichlorobenzene-d4 | 107 | | | 70-13 | 0 % | | " | " | | | " | |
| 460-00-4 | % Bromofluorobenzene | 91 | | | 70-13 | 0 % | | " | " | | | " | |
| 1868-53-7 | % Dibromofluoromethane | 108 | | | 70-13 | 0 % | | " | " | " | | " | |
| 2037-26-5 | % Toluene-d8 | 94 | | | 70-13 | 0 % | | " | " | " | " | " | |
| Analysis pe | erformed by Phoenix Environi | nental Labs, | Inc. * - MAC | T007 | | | | | | | | | |
| 64-17-5 | Ethanol | < 400 | | ug/L | 400 | 400 | 1 | SW8260C. | " | 09-Jan-19 10:09 | M-CT007 | 462669A | ı |
| | acted Analyses by method SW3520C | | | | | | | | | | | | |
| Analysis pe | erformed by Phoenix Environi | nental Labs, | Inc. * - MAC | T007 | | | | | | | | | |
| 50-32-8 | Benzo(a)pyrene | < 0.10 | | ug/L | 0.10 | 0.10 | 1 | SW8270D (SIM) | 22-Dec-18 | 27-Dec-18 02:45 | M-CT007 | 461099A | i |
| 91-20-3 | Naphthalene | < 0.48 | | ug/L | 0.48 | 0.48 | 1 | п | " | II | " | " | |
| Surrogate i | recoveries: | | | | | | | | | | | | |
| 321-60-8 | % 2-Fluorobiphenyl | 52 | | | 30-13 | 0 % | | 11 | " | " | " | " | |
| 4165-60-0 | % Nitrobenzene-d5 | 44 | | | 30-13 | 0 % | | " | " | " | | " | |

30-130 %

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98904-43-9 % Terphenyl-d14

31

| Sample Id Trip Blar SC52773 | | | | | Project # Chelsea | | <u>Matrix</u> Trip Blan | | ection Date 9-Dec-18 00 | | - | eived Dec-18 | |
|-----------------------------------|------------------------------|------------------|-------------|-------|----------------------|------|----------------------------|-------------|-------------------------|--------------------|---------|-----------------|-------|
| CAS No. | Analyte(s) | Result | Flag | Units | *RDL | MDL | Dilution | Method Ref. | Prepared | Analyzed | Analyst | Batch | Cert. |
| Subcontra | acted Analyses | | | | | | | | | | | | |
| Subcontra | acted Analyses | | | | | | | | | | | | |
| <u>Prepared</u> | by method SW8260C | | | | | | | | | | | | |
| Analysis pe | erformed by Phoenix Environr | nental Labs, Ind | c. * - MACT | 7007 | | | | | | | | | |
| 71-43-2 | Benzene | < 0.70 | | ug/L | 0.70 | 0.70 | 1 | SW8260C | 19-Dec-18 | 24-Dec-18 11:19 | M-CT007 | 461279A | |
| 1634-04-4 | Methyl t-butyl ether (MTBE) | < 1.0 | | ug/L | 1.0 | 1.0 | 1 | " | " | " | " | " | |
| 91-20-3 | Naphthalene | < 1.0 | | ug/L | 1.0 | 1.0 | 1 | " | " | " | " | " | |
| Surrogate | recoveries: | | | | | | | | | | | | |
| 2199-69-1 | % 1,2-dichlorobenzene-d4 | 110 | | | 70-13 | 0 % | | " | " | " | " | " | |
| 460-00-4 | % Bromofluorobenzene | 86 | | | 70-13 | 0 % | | " | " | " | " | " | |
| 1868-53-7 | % Dibromofluoromethane | 109 | | | 70-13 | 0 % | | " | " | " | " | " | |
| 2037-26-5 | % Toluene-d8 | 97 | | | 70-13 | 0 % | | u u | " | " | " | " | |
| Analysis pe | erformed by Phoenix Environn | nental Labs, Inc | c. * - MACT | 7007 | | | | | | | | | |
| 64-17-5 | Ethanol | < 400 | | ug/L | 400 | 400 | 1 | SW8260C. | " | 09-Jan-19 09:47 | M-CT007 | 462669A | |

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General Chemistry Parameters - Quality Control

| Analyte(s) | Result | Flag | Units | *RDL | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit |
|-------------------------------------|--------|------|------------|-----------------|----------------|------------------|-------------|----------------|-----|--------------|
| ASTM D 1293-99B | | | | | | | | | | |
| Batch 1816206 - General Preparation | | | | | | | | | | |
| <u>Duplicate (1816206-DUP1)</u> | | | Source: SC | <u>52773-01</u> | Pre | epared & Aı | nalyzed: 21 | I-Dec-18 | | |
| pH | 7.01 | | pH Units | | | 7.03 | | | 0.3 | 5 |
| Reference (1816206-SRM1) | | | | | Pre | epared & Ar | nalyzed: 21 | I-Dec-18 | | |
| рН | 5.99 | | pH Units | | 6.00 | | 100 | 97.5-102. 5 | | |
| Reference (1816206-SRM2) | | | | | Pre | epared & Ar | nalyzed: 21 | I-Dec-18 | | |
| рН | 6.04 | | pH Units | | 6.00 | | 101 | 97.5-102. 5 | | |

| | <u>.</u> . | | | | Spike | Source | 0/=== | %REC | | RPD |
|-----------------------------|------------|------|------------|------|-------|------------|--------------------|---------|-----|-------|
| Analyte(s) | Result | Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limit |
| E1664A | | | | | | | | | | |
| Batch 461365A - E1664A | | | | | | | | | | |
| BLK (CC19445-BLK) | | | | | Pre | epared & A | nalyzed: 27- | -Dec-18 | | |
| Oil and Grease by EPA 1664A | < 1.4 | c3 | mg/L | 1.4 | 40 | | BRL | | | |
| LCS (CC19445-LCS) | | | · · | | Pre | epared: A | nalyzed: 27 | -Dec-18 | | |
| Oil and Grease by EPA 1664A | 38.70 | сЗ | mg/L | 1.4 | 40 | | 97 | 85-115 | | 20 |
| LCSD (CC19445-LCSD) | | | 3 | | | epared: A | nalyzed: 27 | | | |
| Oil and Grease by EPA 1664A | 38.30 | сЗ | mg/L | 1.4 | 40 | <u> </u> | 96 | 85-115 | 1.0 | 20 |
| SM2540D-11 | | | Ü | | | | | | | |
| Batch 461248A - SM2540D-11 | | | | | | | | | | |
| | | | | | Dec | anarad: A | naluzadi 06 | Dog 19 | | |
| BLK (CC18506-BLK) | < 5.0 | | ma/l | 5.0 | 61.3 | epared: A | nalyzed: 26 BRL | -Dec-16 | | |
| Total Suspended Solids | < 5.0 | | mg/L | | | | | | | |
| DUP (CC18506-DUP) | . 5.0 | | Source: CC | | | epared: A | nalyzed: 26 | | NO | |
| Total Suspended Solids | < 5.0 | | mg/L | 5.0 | 61.3 | | | - | NC | |
| LCS (CC18506-LCS) | | | - | | | epared: A | nalyzed: 26 | | | |
| Total Suspended Solids | 54.00 | | mg/L | 5.0 | 61.3 | | 88 | 85-115 | | |
| <u>SW8260C</u> | | | | | | | | | | |
| Batch 461279A - SW8260C | | | | | | | | | | |
| BLK (CC19490-BLK) | | | | | Pre | epared: A | nalyzed: 24 | -Dec-18 | | |
| Carbon tetrachloride | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| cis-1,3-Dichloropropene | ND | c1 | ug/L | 0.40 | | | ND | - | | |
| Hexachlorobutadiene | ND | c1 | ug/L | 0.40 | | | ND | - | | |
| Ethylbenzene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Ethyl ether | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Dichlorodifluoromethane | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Dibromomethane | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Dibromochloromethane | ND | c1 | ug/L | 0.50 | | | ND | - | | |
| Isopropylbenzene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| cis-1,2-Dichloroethene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Chloromethane | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Chloroform | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| m&p-Xylene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Chlorobenzene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| p-lsopropyltoluene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Carbon Disulfide | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Bromomethane | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Chloroethane | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| tert-Butylbenzene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Vinyl chloride | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Trichlorofluoromethane | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Trichloroethene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| trans-1,4-dichloro-2-butene | ND | c1 | ug/L | 5.0 | | | ND | - | | |
| trans-1,3-Dichloropropene | ND | c1 | ug/L | 0.40 | | | ND | - | | |
| trans-1,2-Dichloroethene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Toluene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| n-Propylbenzene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Tetrachloroethene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Methyl ethyl ketone | ND | c1 | ug/L | 5.0 | | | ND | - | | |
| Styrene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| sec-Butylbenzene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Trichlorotrifluoroethane | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| o-Xylene | ND | c1 | ug/L | 1.0 | | | ND | - | | |

| analyte(s) | Result | Flag | Units | *RDL | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit |
|-------------------------------------|----------|------|--------------|------|----------------|------------------|-------------|----------------|-----|--------------|
| W8260C | | | | | | | | | | |
| atch 461279A - SW8260C | | | | | | | | | | |
| BLK (CC19490-BLK) | | | | | Pre | epared: A | nalyzed: 24 | -Dec-18 | | |
| Bromoform | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| n-Butylbenzene | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| Methylene chloride | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| Tetrahydrofuran (THF) | ND | c1 | ug/L | 2.5 | | | ND | _ | | |
| 1,1-Dichloroethene | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| 1,2-Dichlorobenzene | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| 1,2-Dibromoethane | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| 1,2-Dibromo-3-chloropropane | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| 1,2,4-Trimethylbenzene | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| 1,2,4-Trichlorobenzene | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| 1,2,3-Trichloropropane | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| 1,2-Dichloroethane | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| 1,1-Dichloropropene | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| 1,2-Dichloropropane | ND | c1 | ug/L | 1.0 | | | ND | _ | | |
| 1,1-Dichloroethane | ND | c1 | ug/L ug/L | 1.0 | | | ND | - | | |
| 1,1,2-Trichloroethane | ND ND | c1 | | 1.0 | | | ND | - | | |
| | | c1 | ug/L | | | | ND | - | | |
| 1,1,2,2-Tetrachloroethane | ND | c1 | ug/L | 0.50 | | | | | | |
| 1,1,1-Trichloroethane | ND | | ug/L | 1.0 | | | ND | - | | |
| 1,1,1,2-Tetrachloroethane | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Bromodichloromethane | ND | c1 | ug/L | 0.50 | | | ND | - | | |
| 1,2,3-Trichlorobenzene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| 2-Chlorotoluene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Bromochloromethane | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Bromobenzene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Acrylonitrile | ND | c1 | ug/L | 5.0 | | | ND | - | | |
| Acetone | ND | c1 | ug/L | 5.0 | | | ND | - | | |
| 4-Methyl-2-pentanone | ND | c1 | ug/L | 5.0 | | | ND | - | | |
| 4-Chlorotoluene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| 2-Hexanone | ND | c1 | ug/L | 5.0 | | | ND | - | | |
| 2,2-Dichloropropane | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| 1,4-dioxane | ND | c1 | ug/L | 100 | | | ND | - | | |
| 1,4-Dichlorobenzene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| 1,3-Dichloropropane | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| 1,3-Dichlorobenzene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| 1,3,5-Trimethylbenzene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| 2-Isopropyltoluene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Methyl t-butyl ether (MTBE) | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Naphthalene | ND | c1 | ug/L | 1.0 | | | ND | - | | |
| Benzene | ND | c1 | ug/L | 0.70 | | | ND | - | | |
| Surrogate: % 1,2-dichlorobenzene-d4 | 105 | c1 | ug/L | | 10 | | 105 | 70-130 | | |
| Surrogate: % Bromofluorobenzene | 89 | c1 | ug/L | | 10 | | 89 | 70-130 | | |
| Surrogate: % Dibromofluoromethane | 102 | c1 | ug/L | | 10 | | 102 | 70-130 | | |
| Surrogate: % Toluene-d8 | 92 | c1 | ug/L | | 10 | | 92 | 70-130 | | |
| LCS (CC19490-LCS) | 02 | | ~g, ∟ | | | epared: A | nalyzed: 24 | | | |
| | 0.000 | c1 | ua/l | 0.40 | 10 | spaieu. A | 100 | | | 30 |
| cis-1,3-Dichloropropene | 9.962 | c1 | ug/L | 0.40 | | | | 70-130 | | |
| m&p-Xylene | 19.30 | c1 | ug/L | 1.0 | 20 | | 96 | 70-130 | | 30 |
| Isopropylbenzene | 9.291 | c1 | ug/L | 1.0 | 10 | | 93 | 70-130 | | 30 |
| Hexachlorobutadiene | 9.398 | c1 | ug/L | 0.40 | 10 | | 94 | 70-130 | | 30 |
| Ethylbenzene | 9.238 | c1 | ug/L | 1.0 | 10 | | 92 | 70-130 | | 30 |
| Ethyl ether | 11.15 | c1 | ug/L | 1.0 | 10 | | 111 | 70-130 | | 30 |

| 6.447 8.137 10.01 9.411 9.488 7.210 9.165 8.488 8.889 8.834 | c1 c1 c1 c1 c1 c1 c1 c1 | ug/L ug/L ug/L ug/L ug/L ug/L | 1.0 1.0 0.50 1.0 | <u>Pro</u> 10 10 10 | epared: A | <u>nalyzed: 24</u> 64 81 | <u>-Dec-18</u> 40-160 70-130 | | 30 |
|--|---|---|---|--|---|--|---|---|--|
| 8.137 10.01 9.411 9.488 7.210 9.165 8.488 8.889 | c1 c1 c1 c1 c1 c1 | ug/L ug/L ug/L ug/L | 1.0 0.50 | 10 10 | epared: A | 64 81 | 40-160 | | 30 |
| 8.137 10.01 9.411 9.488 7.210 9.165 8.488 8.889 | c1 c1 c1 c1 c1 c1 | ug/L ug/L ug/L ug/L | 1.0 0.50 | 10 10 | epared: A | 64 81 | 40-160 | | 30 |
| 8.137 10.01 9.411 9.488 7.210 9.165 8.488 8.889 | c1 c1 c1 c1 c1 c1 | ug/L ug/L ug/L ug/L | 1.0 0.50 | 10 | | 81 | | | 30 |
| 10.01 9.411 9.488 7.210 9.165 8.488 8.889 | c1 c1 c1 c1 c1 | ug/L ug/L ug/L ug/L | 0.50 | | | | 70-130 | | |
| 9.411 9.488 7.210 9.165 8.488 8.889 | c1 c1 c1 c1 | ug/L ug/L | | 10 | | | | | 30 |
| 9.488 7.210 9.165 8.488 8.889 | c1 c1 c1 | ug/L ug/L | | | | 100 | 70-130 | | 30 |
| 9.488 7.210 9.165 8.488 8.889 | c1 c1 | ug/L | | 10 | | 94 | 70-130 | | 30 |
| 7.210 9.165 8.488 8.889 | c1 | - | 1.0 | 10 | | 95 | 70-130 | | 30 |
| 9.165 8.488 8.889 | | | 1.0 | 10 | | 72 | 40-160 | | 30 |
| 8.488 8.889 | | ug/L | 1.0 | 10 | | 92 | 70-130 | | 30 |
| 8.889 | | ug/L | 1.0 | 10 | | 85 | 70-130 | | 30 |
| | c1 | ug/L | 1.0 | 10 | | 89 | 70-130 | | 30 |
| 0.004 | c1 | ug/L | 1.0 | 10 | | 88 | 70-130 | | 30 |
| 8.606 | c1 | ug/L | 1.0 | 10 | | 86 | 70-130 | | 30 |
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| | | ug/L | | | | | | | 30 |
| 25.38 | c1 | ug/L | 2.5 | 25 | | 102 | 70-130 | | 30 |
| 9.720 | c1 | ug/L | 1.0 | 10 | | 97 | 70-130 | | 30 |
| 9.582 | c1 | ug/L | 1.0 | 10 | | 96 | 70-130 | | 30 |
| 9.930 | c1 | ug/L | 1.0 | 10 | | 99 | 70-130 | | 30 |
| 10.02 | c1 | ug/L | 1.0 | 10 | | 100 | 70-130 | | 30 |
| 9.667 | c1 | ug/L | 1.0 | 10 | | 97 | 70-130 | | 30 |
| 9.817 | c1 | ug/L | 1.0 | 10 | | 98 | 70-130 | | 30 |
| 9.377 | c1 | ug/L | 1.0 | 10 | | 94 | 70-130 | | 30 |
| 9.197 | c1 | ug/L | 1.0 | 10 | | 92 | 70-130 | | 30 |
| 9.065 | c1 | ug/L | 1.0 | 10 | | 91 | 70-130 | | 30 |
| 8.840 | c1 | ug/L | 1.0 | 10 | | 88 | 70-130 | | 30 |
| 9.139 | c1 | ug/L | 1.0 | 10 | | 91 | 70-130 | | 30 |
| 9.129 | c1 | ug/L | 1.0 | 10 | | 91 | 70-130 | | 30 |
| 9.910 | c1 | ug/L | 1.0 | 10 | | 99 | 70-130 | | 30 |
| 9.707 | c1 | ug/L | 1.0 | 10 | | 97 | 70-130 | | 30 |
| 9.078 | c1 | ug/L | 1.0 | 10 | | 91 | 70-130 | | 30 |
| 9.383 | c1 | - | 1.0 | 10 | | 94 | 70-130 | | 30 |
| | c1 | - | | 10 | | 93 | 70-130 | | 30 |
| | c1 | - | | | | | | | 30 |
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| | 9.582 9.930 10.02 9.667 9.817 9.377 9.197 9.065 8.840 9.139 9.129 9.910 9.707 | 8.497 c1 6.333 c1 8.971 c1 8.280 c1 8.929 c1 54.35 c1 9.602 c1 9.373 c1 10.76 c1 25.38 c1 9.720 c1 9.582 c1 9.930 c1 10.02 c1 9.667 c1 9.817 c1 9.377 c1 9.197 c1 9.065 c1 8.840 c1 9.139 c1 9.129 c1 9.707 c1 9.707 c1 9.707 c1 9.707 c1 9.707 c1 9.707 c1 9.708 c1 9.707 c1 9.707 c1 9.707 c1 9.708 c1 9.383 c1 9.270 c1 9.746 c1 9.515 c1 8.829 c1 9.301 c1 9.498 c1 9.498 c1 9.498 c1 9.304 c1 9.304 c1 9.304 c1 9.304 c1 | 8.497 c1 ug/L 6.333 c1 ug/L 8.971 c1 ug/L 8.280 c1 ug/L 8.929 c1 ug/L 9.602 c1 ug/L 9.373 c1 ug/L 25.38 c1 ug/L 9.720 c1 ug/L 9.582 c1 ug/L 9.930 c1 ug/L 9.930 c1 ug/L 9.930 c1 ug/L 9.377 c1 ug/L 9.197 c1 ug/L 9.199 c1 ug/L 9.199 c1 ug/L 9.199 c1 ug/L 9.129 c1 ug/L 9.270 c1 ug/L 9.383 c1 ug/L 9.395 c1 ug/L 9.301 c1 ug/L 9.301 c1 ug/L 9.301 c1 ug/L 9.304 c1 ug/L 9.304 c1 ug/L 9.304 c1 ug/L | 8.497 c1 ug/L 1.0 6.333 c1 ug/L 1.0 8.971 c1 ug/L 1.0 8.280 c1 ug/L 1.0 8.929 c1 ug/L 1.0 54.35 c1 ug/L 5.0 9.602 c1 ug/L 0.40 9.373 c1 ug/L 1.0 10.76 c1 ug/L 5.0 25.38 c1 ug/L 1.0 9.582 c1 ug/L 1.0 9.582 c1 ug/L 1.0 9.930 c1 ug/L 1.0 9.930 c1 ug/L 1.0 9.9407 c1 ug/L 1.0 9.667 c1 ug/L 1.0 9.377 c1 ug/L 1.0 9.377 c1 ug/L 1.0 9.377 c1 ug/L 1.0 9.197 c1 ug/L 1.0 9.197 c1 ug/L 1.0 9.197 c1 ug/L 1.0 9.105 c1 ug/L 1.0 9.105 c1 ug/L 1.0 9.1065 c1 ug/L 1.0 9.107 c1 ug/L 1.0 9.108 c1 ug/L 1.0 9.109 c1 ug/L 1.0 | 8.497 c1 ug/L 1.0 10 6.333 c1 ug/L 1.0 10 8.971 c1 ug/L 1.0 10 8.280 c1 ug/L 1.0 10 8.929 c1 ug/L 1.0 10 54.35 c1 ug/L 5.0 50 9.602 c1 ug/L 0.40 10 9.373 c1 ug/L 5.0 10 10.76 c1 ug/L 5.0 10 25.38 c1 ug/L 2.5 25 9.720 c1 ug/L 1.0 10 9.582 c1 ug/L 1.0 10 9.582 c1 ug/L 1.0 10 10.02 c1 ug/L 1.0 10 10.02 c1 ug/L 1.0 10 9.667 c1 ug/L 1.0 10 9.667 c1 ug/L 1.0 10 9.377 c1 ug/L 1.0 10 9.390 c1 ug/L 1.0 10 9.391 c1 ug/L 1.0 10 9.391 c1 ug/L 1.0 10 9.397 c1 ug/L 1.0 10 9.197 c1 ug/L 1.0 10 9.197 c1 ug/L 1.0 10 9.198 c1 ug/L 1.0 10 9.109 c1 ug/L 1.0 10 9.119 c1 ug/L 1.0 10 9.129 c1 ug/L 1.0 10 9.707 c1 ug/L 1.0 10 9.708 c1 ug/L 1.0 10 9.709 c1 ug/L 1.0 10 9.700 c1 ug/L 1.0 10 | 8.497 c1 ug/L 1.0 10 6.333 c1 ug/L 1.0 10 8.971 c1 ug/L 1.0 10 8.280 c1 ug/L 1.0 10 8.929 c1 ug/L 1.0 10 54.35 c1 ug/L 5.0 50 9.602 c1 ug/L 1.0 10 9.373 c1 ug/L 1.0 10 10.76 c1 ug/L 5.0 10 25.38 c1 ug/L 1.0 10 9.582 c1 ug/L 1.0 10 9.583 c1 ug/L 1.0 10 9.584 c1 ug/L 1 | 8.497 c1 ug/L 1.0 10 85 6.333 c1 ug/L 1.0 10 63 8.971 c1 ug/L 1.0 10 90 8.280 c1 ug/L 1.0 10 83 8.929 c1 ug/L 5.0 50 109 9.602 c1 ug/L 0.40 10 96 9.373 c1 ug/L 5.0 10 10 94 10.76 c1 ug/L 5.0 10 10 97 9.582 c1 ug/L 1.0 10 97 9.582 c1 ug/L 1.0 10 96 9.330 c1 ug/L 1.0 10 96 9.377 c1 ug/L 1.0 10 96 9.377 c1 ug/L 1.0 10 97 9.817 c1 ug/L 1.0 10 99 9.817 c1 ug/L 1.0 10 97 9.817 c1 ug/L 1.0 10 97 9.817 c1 ug/L 1.0 10 97 9.917 c1 ug/L 1.0 10 99 9.377 c1 ug/L 1.0 10 99 9.377 c1 ug/L 1.0 10 99 9.377 c1 ug/L 1.0 10 97 9.817 c1 ug/L 1.0 10 99 9.377 c1 ug/L 1.0 10 99 9.379 c1 ug/L 1.0 10 99 9.379 c1 ug/L 1.0 10 99 9.390 c1 ug/L 1.0 10 99 9.391 c1 ug/L 1.0 10 99 9.391 c1 ug/L 1.0 10 99 9.393 c1 ug/L 1.0 10 99 9.394 c1 ug/L 1.0 10 99 9.395 c1 ug/L 1.0 10 99 9.394 c1 ug/L 1.0 10 99 | 8.497 c1 ug/L 1.0 10 85 70-130 6.333 c1 ug/L 1.0 10 63 40-160 8.971 c1 ug/L 1.0 10 90 70-130 8.280 c1 ug/L 1.0 10 83 70-130 8.929 c1 ug/L 5.0 50 109 70-130 9.602 c1 ug/L 0.40 10 96 70-130 9.602 c1 ug/L 5.0 10 108 40-160 25.38 c1 ug/L 5.0 10 108 40-160 25.38 c1 ug/L 1.0 10 97 70-130 9.720 c1 ug/L 1.0 10 97 70-130 9.822 c1 ug/L 1.0 10 96 70-130 9.330 c1 ug/L 1.0 10 97 70-130 | 8.497 c1 ug/L 1.0 10 85 70-130 6.333 c1 ug/L 1.0 10 63 40-160 8.971 c1 ug/L 1.0 10 90 70-130 8.280 c1 ug/L 1.0 10 83 70-130 8.929 c1 ug/L 1.0 10 89 70-130 9.602 c1 ug/L 5.0 50 109 70-130 9.373 c1 ug/L 1.0 10 96 70-130 9.373 c1 ug/L 5.0 10 10 94 70-130 10.76 c1 ug/L 5.0 10 108 40-160 25.38 c1 ug/L 2.5 25 102 70-130 9.582 c1 ug/L 1.0 10 97 70-130 9.582 c1 ug/L 1.0 10 99 70-130 9.582 c1 ug/L 1.0 10 99 70-130 9.817 c1 ug/L 1.0 10 99 70-130 9.667 c1 ug/L 1.0 10 99 70-130 9.670 c1 ug/L 1.0 10 99 70-130 9.377 c1 ug/L 1.0 10 99 70-130 9.687 c1 ug/L 1.0 10 99 70-130 9.687 c1 ug/L 1.0 10 99 70-130 9.698 c1 ug/L 1.0 10 99 70-130 9.197 c1 ug/L 1.0 10 99 70-130 9.199 c1 ug/L 1.0 10 91 70-130 9.190 c1 ug/L 1.0 10 91 70-130 9.190 c1 ug/L 1.0 10 91 70-130 9.190 c1 ug/L 1.0 10 99 70-130 9.190 c1 ug/L 1.0 10 99 70-130 9.190 c1 ug/L 1.0 10 99 70-130 9.370 c1 ug/L 1.0 10 99 70-130 9.370 c1 ug/L 1.0 10 99 70-130 9.383 c1 ug/L 1.0 10 99 70-130 9.391 c1 ug/L 1.0 10 99 70-130 9.3920 c1 ug/L 1.0 10 99 70-130 9.3930 c1 ug/L 1.0 10 99 70-130 9.3948 c1 ug/L 1.0 10 99 70-130 9.3948 c1 ug/L 1.0 10 99 70-130 9.3949 c1 ug/L 1.0 10 99 70-130 9.3940 c1 ug/L 1.0 10 99 70-130 |

| Analyte(s) | Result | Flag | Units | *RDL | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit |
|-------------------------------------|--------|------|--------------|------|----------------|------------------|-------------|------------------|------|--------------|
| SW8260C | | | | | | | | | | |
| Batch 461279A - SW8260C | | | | | | | | | | |
| LCS (CC19490-LCS) | | | | | Pr | epared: Ar | nalyzed: 24 | -Dec-18 | | |
| Bromochloromethane | 9.789 | c1 | ug/L | 1.0 | 10 | | 98 | 70-130 | | 30 |
| 1,2,4-Trimethylbenzene | 9.841 | c1 | ug/L | 1.0 | 10 | | 98 | 70-130 | | 30 |
| Bromobenzene | 9.374 | c1 | ug/L | 1.0 | 10 | | 94 | 70-130 | | 30 |
| 1,3,5-Trimethylbenzene | 9.418 | c1 | ug/L | 1.0 | 10 | | 94 | 70-130 | | 30 |
| Acetone | 9.175 | c1 | ug/L | 5.0 | 10 | | 92 | 40-160 | | 30 |
| 4-Methyl-2-pentanone | 10.36 | c1 | ug/L | 5.0 | 10 | | 104 | 40-160 | | 30 |
| 4-Chlorotoluene | 9.233 | c1 | ug/L | 1.0 | 10 | | 92 | 70-130 | | 30 |
| 2-Hexanone | 9.898 | c1 | ug/L | 5.0 | 10 | | 99 | 40-160 | | 30 |
| 2-Chlorotoluene | 9.324 | c1 | ug/L | 1.0 | 10 | | 93 | 70-130 | | 30 |
| 2,2-Dichloropropane | 8.928 | c1 | ug/L | 1.0 | 10 | | 89 | 70-130 | | 30 |
| 1,4-dioxane | 177.1 | c1 | ug/L | 100 | 200 | | 89 | 40-160 | | 30 |
| 1,3-Dichloropropane | 9.644 | c1 | ug/L | 1.0 | 10 | | 96 | 70-130 | | 30 |
| 2-Isopropyltoluene | 10.01 | c1 | ug/L | 1.0 | 10 | | 100 | 70-130 | | 30 |
| Naphthalene | 10.62 | c1 | ug/L | 1.0 | 10 | | 106 | 70-130 | | 30 |
| Benzene | 9.012 | c1 | ug/L ug/L | 0.70 | 10 | | 90 | 70-130 | | 30 |
| Methyl t-butyl ether (MTBE) | 9.504 | c1 | ug/L | 1.0 | 10 | | 95 | 70-130 | | 30 |
| Surrogate: % Toluene-d8 | 10.21 | c1 | ug/L | | 10 | | 102 | 70-130 | | |
| Surrogate: % Dibromofluoromethane | 10.65 | c1 | ug/L | | 10 | | 106 | 70-130 | | |
| Surrogate: % 1,2-dichlorobenzene-d4 | 10.10 | c1 | ug/L | | 10 | | 101 | 70-130 70-130 | | |
| Surrogate: % Bromofluorobenzene | 10.35 | c1 | ug/L | | 10 | | 104 | 70-130 70-130 | | |
| • | 70.00 | - | ug/L | | | epared: Ar | nalyzed: 24 | | | |
| LCSD (CC19490-LCSD) | 9.690 | c1 | ua/l | 1.0 | 10 | epareu. Ai | 97 | 70-130 | 0.0 | 30 |
| 1,2,3-Trichlorobenzene | | c1 | ug/L | | 10 | | | | | 30 |
| 1,1,1,2-Tetrachloroethane | 9.385 | c1 | ug/L | 1.0 | | | 94 | 70-130 | 1.1 | |
| 1,1,1-Trichloroethane | 9.403 | | ug/L | 1.0 | 10 | | 94 | 70-130 | 4.3 | 30 |
| 1,1,2,2-Tetrachloroethane | 9.843 | c1 | ug/L | 0.50 | 10 | | 98 | 70-130 | 1.0 | 30 |
| 1,1,2-Trichloroethane | 10.02 | c1 | ug/L | 1.0 | 10 | | 100 | 70-130 | 5.1 | 30 |
| 1,1-Dichloropropene | 9.822 | c1 | ug/L | 1.0 | 10 | | 98 | 70-130 | 10.8 | 30 |
| 1,1-Dichloroethene | 9.422 | c1 | ug/L | 1.0 | 10 | | 94 | 70-130 | 6.6 | 30 |
| Ethyl ether | 10.84 | c1 | ug/L | 1.0 | 10 | | 108 | 70-130 | 2.7 | 30 |
| Chlorobenzene | 9.372 | c1 | ug/L | 1.0 | 10 | | 94 | 70-130 | 5.5 | 30 |
| n-Butylbenzene | 9.749 | c1 | ug/L | 1.0 | 10 | | 97 | 70-130 | 3.1 | 30 |
| Methylene chloride | 9.795 | c1 | ug/L | 1.0 | 10 | | 98 | 70-130 | 1.0 | 30 |
| Methyl ethyl ketone | 10.82 | c1 | ug/L | 5.0 | 10 | | 108 | 40-160 | 0.0 | 30 |
| m&p-Xylene | 19.62 | c1 | ug/L | 1.0 | 20 | | 98 | 70-130 | 2.1 | 30 |
| Isopropylbenzene | 9.845 | c1 | ug/L | 1.0 | 10 | | 98 | 70-130 | 5.2 | 30 |
| o-Xylene | 9.863 | c1 | ug/L | 1.0 | 10 | | 99 | 70-130 | 1.0 | 30 |
| Ethylbenzene | 9.526 | c1 | ug/L | 1.0 | 10 | | 95 | 70-130 | 3.2 | 30 |
| p-Isopropyltoluene | 10.16 | c1 | ug/L | 1.0 | 10 | | 102 | 70-130 | 5.0 | 30 |
| Dibromomethane | 10.27 | c1 | ug/L | 1.0 | 10 | | 103 | 70-130 | 5.0 | 30 |
| cis-1,3-Dichloropropene | 10.53 | c1 | ug/L | 0.40 | 10 | | 105 | 70-130 | 4.9 | 30 |
| cis-1,2-Dichloroethene | 9.486 | c1 | ug/L | 1.0 | 10 | | 95 | 70-130 | 0.0 | 30 |
| Chloromethane | 7.683 | c1 | ug/L | 1.0 | 10 | | 77 | 40-160 | 6.7 | 30 |
| Chloroform | 9.324 | c1 | ug/L | 1.0 | 10 | | 93 | 70-130 | 1.1 | 30 |
| Chloroethane | 9.066 | c1 | ug/L | 1.0 | 10 | | 91 | 70-130 | 6.8 | 30 |
| Hexachlorobutadiene | 9.185 | c1 | ug/L | 0.40 | 10 | | 92 | 70-130 | 2.2 | 30 |
| trans-1,2-Dichloroethene | 9.538 | c1 | ug/L | 1.0 | 10 | | 95 | 70-130 | 1.1 | 30 |
| 1,2,3-Trichloropropane | 9.364 | c1 | ug/L | 1.0 | 10 | | 94 | 70-130 | 1.1 | 30 |
| 1,1-Dichloroethane | 9.404 | c1 | ug/L | 1.0 | 10 | | 94 | 70-130 | 1.1 | 30 |
| Vinyl chloride | 8.734 | c1 | ug/L | 1.0 | 10 | | 87 | 70-130 | 7.1 | 30 |
| Trichlorotrifluoroethane | 10.01 | c1 | ug/L | 1.0 | 10 | | 100 | 70-130 | 10.5 | 30 |

| nalyte(s) | Result | Flag | Units | *RDL | Spike Level | Source Result | %REC | %REC Limits | RPD | RPI Lim |
|---|----------------|----------|-------|------------|----------------|------------------|-------------|------------------|------------|------------|
| W8260C | | | | | | | | | | |
| atch 461279A - SW8260C | | | | | | | | | | |
| LCSD (CC19490-LCSD) | | | | | Pre | epared: A | nalyzed: 24 | -Dec-18 | | |
| Trichlorofluoromethane | 9.035 | c1 | ug/L | 1.0 | 10 | , | 90 | 70-130 | 8.1 | 30 |
| Trichloroethene | 9.490 | c1 | ug/L | 1.0 | 10 | | 95 | 70-130 | 6.5 | 30 |
| n-Propylbenzene | 9.732 | c1 | ug/L | 1.0 | 10 | | 97 | 70-130 | 3.1 | 30 |
| trans-1,3-Dichloropropene | 9.918 | c1 | ug/L | 0.40 | 10 | | 99 | 70-130 | 3.1 | 30 |
| Dibromochloromethane | 10.08 | c1 | ug/L | 0.50 | 10 | | 101 | 70-130 | 1.0 | 30 |
| Toluene | 9.931 | c1 | ug/L | 1.0 | 10 | | 99 | 70-130 | 8.4 | 30 |
| Tetrahydrofuran (THF) | 25.22 | c1 | ug/L | 2.5 | 25 | | 101 | 70-130 | 1.0 | 30 |
| Tetrachloroethene | 9.892 | c1 | ug/L | 1.0 | 10 | | 99 | 70-130 | 15.2 | 30 |
| tert-Butylbenzene | 9.971 | c1 | ug/L | 1.0 | 10 | | 100 | 70-130 | 4.1 | 30 |
| Styrene | 10.05 | c1 | ug/L | 1.0 | 10 | | 100 | 70-130 | 1.0 | 30 |
| sec-Butylbenzene | 10.44 | c1 | ug/L | 1.0 | 10 | | 104 | 70-130 | 3.9 | 30 |
| trans-1,4-dichloro-2-butene | 55.53 | c1 | ug/L | 5.0 | 50 | | 111 | 70-130 | 1.8 | 30 |
| 1,2-Dichloropropane | 10.39 | c1 | ug/L | 1.0 | 10 | | 104 | 70-130 | 11.2 | 30 |
| 2,2-Dichloropropane | 9.379 | c1 | ug/L | 1.0 | 10 | | 94 | 70-130 | 5.5 | 30 |
| 1,4-dioxane | 197.2 | c1 | ug/L | 100 | 200 | | 99 | 40-160 | 10.6 | 30 |
| 1,4-Dichlorobenzene | 9.232 | c1 | ug/L | 1.0 | 10 | | 92 | 70-130 | 1.1 | 30 |
| 1,3-Dichloropropane | 9.496 | c1 | ug/L | 1.0 | 10 | | 95 | 70-130 | 1.0 | 30 |
| Carbon tetrachloride | 9.098 | c1 | ug/L | 1.0 | 10 | | 91 | 70-130 | 3.4 | 30 |
| 2-Chlorotoluene | 9.718 | c1 | ug/L | 1.0 | 10 | | 97 | 70-130 | 4.2 | 30 |
| Dichlorodifluoromethane | 7.098 | c1 | ug/L | 1.0 | 10 | | 71 | 40-160 | 10.4 | 30 |
| 1,3-Dichlorobenzene | 9.470 | c1 | ug/L | 1.0 | 10 | | 95 | 70-130 | 3.2 | 30 |
| 1,2-Dichloroethane | 9.578 | c1 | ug/L | 1.0 | 10 | | 96 | 70-130 | 5.2 | 30 |
| 1,2-Dichlorobenzene | 9.370 | c1 | ug/L | 1.0 | 10 | | 94 | 70-130 | 3.2 | 30 |
| 1,2-Dibromoethane | 9.559 | c1 | ug/L | 1.0 | 10 | | 96 | 70-130 | 1.0 | 30 |
| | | c1 | - | | 10 | | 104 | 70-130 | | 30 |
| 1,2-Dibromo-3-chloropropane | 10.39 10.07 | c1 | ug/L | 1.0 1.0 | 10 | | 104 | 70-130 70-130 | 7.0 3.0 | 30 |
| 1,2,4-Trimethylbenzene 1,2,4-Trichlorobenzene | 9.544 | c1 | ug/L | 1.0 | 10 | | 95 | 70-130 | | 30 |
| 1,3,5-Trimethylbenzene | 9.968 | c1 | ug/L | 1.0 | 10 | | 100 | 70-130 | 1.1 6.2 | 30 |
| • | | c1 | ug/L | | | | | | | |
| Bromochloromethane Carbon Disulfide | 9.563 | c1 | ug/L | 1.0 1.0 | 10 10 | | 96 90 | 70-130 70-130 | 2.1 4.5 | 30 30 |
| | 8.972 | | ug/L | | | | | | | |
| Bromomethane Bromodichloromethane | 6.757 | c1 | ug/L | 1.0 | 10 | | 68 | 40-160 | 7.6 | 30 |
| | 9.976 | c1 c1 | ug/L | 0.50 | 10 | | 100 | 70-130 | 6.2 | 30 |
| 2-Hexanone | 10.42 | | ug/L | 5.0 | 10 | | 104 | 40-160 | 4.9 | 30 |
| Bromobenzene | 9.545 | c1 | ug/L | 1.0 | 10 | | 95 | 70-130 | 1.1 | 30 |
| Acrylonitrile | 10.14 | c1 | ug/L | 5.0 | 10 | | 101 | 70-130 | 2.0 | 30 |
| 4-Methyl-2-pentanone | 10.71 | c1 | ug/L | 5.0 | 10 | | 107 | 40-160 | 2.8 | 30 |
| Bromoform | 10.43 | c1 | ug/L | 1.0 | 10 | | 104 | 70-130 | 4.9 | 30 |
| 4-Chlorotoluene | 9.517 | c1 | ug/L | 1.0 | 10 | | 95 106 | 70-130 | 3.2 | 30 |
| 2-Isopropyltoluene | 10.56 | c1 | ug/L | 1.0 | 10 | | 106 | 70-130 | 5.8 | 30 |
| Acetone | 9.739 | c1 | ug/L | 5.0 | 10 | | 97 | 40-160 | 5.3 | 30 |
| Benzene | 9.563 | c1 | ug/L | 0.70 | 10 | | 96 | 70-130 | 6.5 | 30 |
| Naphthalene | 10.75 | c1 | ug/L | 1.0 | 10 | | 108 | 70-130 | 1.9 | 30 |
| Methyl t-butyl ether (MTBE) | 9.611 | c1 | ug/L | 1.0 | 10 | | 96 | 70-130 | 1.0 | 30 |
| Surrogate: % Bromofluorobenzene | 9.919 | c1 | ug/L | | 10 | | 99 | 70-130 | | |
| Surrogate: % Dibromofluoromethane | 10.24 | c1 | ug/L | | 10 | | 102 | 70-130 | | |
| Surrogate: % Toluene-d8 | 10.39 | c1 | ug/L | | 10 | | 104 | 70-130 | | |
| Surrogate: % 1,2-dichlorobenzene-d4 | 10.03 | c1 | ug/L | | 10 | | 100 | 70-130 | | |

BLK (CC19490-BLK) Prepared: Analyzed: 24-Dec-18

| | | | | | Spike | Source | | %REC | | RPD |
|-------------------------------|--------|------|------------|------|-----------|------------|----------------------------|------------------|-------|------|
| Analyte(s) | Result | Flag | Units | *RDL | Level | Result | %REC | Limits | RPD | Limi |
| SW8260C. | | | | | | | | | | |
| Batch 462669A - SW8260C | | | | | | | | | | |
| BLK (CC19490-BLK) | | | | | Pro | epared: / | Analyzed: 24 | -Dec-18 | | |
| Ethanol | ND | c4 | ug/L | 200 | <u></u> | <u>.</u> | ND | - | | |
| LCS (CC19490-LCS) | | | -3 | | Pr | epared: / | Analyzed: 24 | -Dec-18 | | |
| Ethanol | 213.4 | c4 | ug/L | 200 | 250 | cparca. 7 | 85 | 70-130 | | 30 |
| LCSD (CC19490-LCSD) | 210.4 | | ug/L | 200 | | oparod: | Analyzed: 24 | | | 00 |
| Ethanol | 197.7 | c4 | ug/L | 200 | 250 | epared: / | 4 <u>naiyzed. 24</u> 79 | 70-130 | 7.3 | 30 |
| | 197.7 | C4 | · · | | | | | | 7.5 | 30 |
| MS (CC19490-MS) | 007.0 | 04 | Source: SO | | | • | Analyzed: 09 | | | 20 |
| Ethanol | 237.8 | c4 | ug/L | 200 | 250 | BRL | 95 | 70-130 | | 30 |
| MSD (CC19490-MSD) | | | Source: SC | | | | Analyzed: 09 | | | |
| Ethanol | 253.0 | c4 | ug/L | 200 | 250 | BRL | 101 | 70-130 | 6.1 | 30 |
| SW8270D (SIM) | | | | | | | | | | |
| Batch 461099A - SW3520C | | | | | | | | | | |
| BLK (CC17546-BLK) | | | | | <u>Pr</u> | epared: 22 | 2-Dec-18 Ar | nalyzed: 26-D | ec-18 | |
| Chrysene | ND | | ug/L | 0.50 | | | ND | - | | |
| Phenanthrene | ND | | ug/L | 0.07 | | | ND | - | | |
| Indeno(1,2,3-cd)pyrene | ND | | ug/L | 0.10 | | | ND | - | | |
| Fluorene | ND | | ug/L | 0.50 | | | ND | - | | |
| Pyrene | ND | | ug/L | 0.50 | | | ND | - | | |
| Dibenz(a,h)anthracene | ND | | ug/L | 0.10 | | | ND | - | | |
| Benzo(b)fluoranthene | ND | | ug/L | 0.08 | | | ND | - | | |
| Benzo(k)fluoranthene | ND | | ug/L | 0.30 | | | ND | - | | |
| Benzo(ghi)perylene | ND | | ug/L | 0.48 | | | ND | - | | |
| Fluoranthene | ND | | ug/L | 0.50 | | | ND | - | | |
| Benz(a)anthracene | ND | | ug/L | 0.06 | | | ND | - | | |
| Anthracene | ND | | ug/L | 0.50 | | | ND | - | | |
| Acenaphthylene | ND | | ug/L | 0.30 | | | ND | _ | | |
| Acenaphthene | ND | | ug/L | 0.50 | | | ND | _ | | |
| 2-Methylnaphthalene | ND | | ug/L | 0.50 | | | ND | _ | | |
| Benzo(a)pyrene | ND | | ug/L | 0.20 | | | ND | _ | | |
| Naphthalene | ND | | ug/L | 0.50 | | | ND | _ | | |
| Surrogate: % Nitrobenzene-d5 | 10 | | | | 5 | | 10 | 30-130 | | |
| = | 76 | | ug/L | | | | | | | |
| Surrogate: % Terphenyl-d14 | 29 | | ug/L | | 5 5 | | 76 29 | 30-130 30-130 | | |
| Surrogate: % 2-Fluorobiphenyl | 29 | | ug/L | | | | | | | |
| LCS (CC17546-LCS) | | | | 0.00 | | epared: 22 | | nalyzed: 26-D | ec-18 | |
| Benzo(b)fluoranthene | 6.315 | | ug/L | 0.08 | 10 | | 63 | 30-130 | | 20 |
| Benzo(k)fluoranthene | 6.628 | | ug/L | 0.30 | 10 | | 66 | 30-130 | | 20 |
| Pyrene | 6.292 | | ug/L | 0.50 | 10 | | 63 | 30-130 | | 20 |
| Phenanthrene | 6.216 | | ug/L | 0.07 | 10 | | 62 | 30-130 | | 20 |
| Indeno(1,2,3-cd)pyrene | 6.352 | | ug/L | 0.10 | 10 | | 64 | 30-130 | | 20 |
| Fluorene | 5.742 | | ug/L | 0.50 | 10 | | 57 | 30-130 | | 20 |
| Fluoranthene | 6.368 | | ug/L | 0.50 | 10 | | 64 | 30-130 | | 20 |
| Dibenz(a,h)anthracene | 7.436 | | ug/L | 0.10 | 10 | | 74 | 30-130 | | 20 |
| Chrysene | 5.998 | | ug/L | 0.50 | 10 | | 60 | 30-130 | | 20 |
| Benz(a)anthracene | 5.319 | | ug/L | 0.06 | 10 | | 53 | 30-130 | | 20 |
| Anthracene | 6.099 | | ug/L | 0.50 | 10 | | 61 | 30-130 | | 20 |
| Acenaphthylene | 4.296 | | ug/L | 0.30 | 10 | | 43 | 30-130 | | 20 |
| Acenaphthene | 5.210 | | ug/L | 0.50 | 10 | | 52 | 30-130 | | 20 |
| 2-Methylnaphthalene | 3.523 | | ug/L | 0.50 | 10 | | 35 | 30-130 | | 20 |
| Benzo(ghi)perylene | 5.874 | | ug/L | 0.48 | 10 | | 59 | 30-130 | | 20 |
| Benzo(a)pyrene | 5.373 | | ug/L | 0.20 | 10 | | 54 | 30-130 | | 20 |

| Analyte(s) | Result | Flag | Units | *RDL | Spike Level | Source Result %I | REC | %REC Limits | RPD | RPE Limi |
|-------------------------------|--------|------|------------|--------|----------------|---------------------|----------------|----------------|-------|-------------|
| SW8270D (SIM) | | | | | | | | | | |
| Batch 461099A - SW3520C | | | | | | | | | | |
| LCS (CC17546-LCS) | | | | | Pre | epared: 22-Dec- | l8 An | alyzed: 26-D | ec-18 | |
| Naphthalene | 2.777 | l, r | ug/L | 0.50 | 10 | 2 | 28 | 30-130 | | 20 |
| Surrogate: % Terphenyl-d14 | 3.913 | | ug/L | | 5 | | 78 | 30-130 | | |
| Surrogate: % 2-Fluorobiphenyl | 1.995 | | ug/L | | 5 | | 10 | 30-130 | | |
| Surrogate: % Nitrobenzene-d5 | 0.9079 | l, r | ug/L | | 5 | | 18 | 30-130 | | |
| LCSD (CC17546-LCSD) | | | Ü | | Pre | epared: 22-Dec- | 18 An: | alvzed: 26-D | ec-18 | |
| Dibenz(a,h)anthracene | 7.372 | | ug/L | 0.10 | 10 | • | 74 | 30-130 | 0.0 | 20 |
| Indeno(1,2,3-cd)pyrene | 6.223 | | ug/L | 0.10 | 10 | (| 62 | 30-130 | 3.2 | 20 |
| Benzo(b)fluoranthene | 6.038 | | ug/L | 0.08 | 10 | | 60 | 30-130 | 4.9 | 20 |
| Fluoranthene | 6.269 | | ug/L | 0.50 | 10 | (| 3 | 30-130 | 1.6 | 20 |
| Pyrene | 6.686 | | ug/L | 0.50 | 10 | (| 67 | 30-130 | 6.2 | 20 |
| Chrysene | 6.100 | | ug/L | 0.50 | 10 | (| 31 | 30-130 | 1.7 | 20 |
| Benzo(k)fluoranthene | 6.882 | | ug/L | 0.30 | 10 | (| 3 9 | 30-130 | 4.4 | 20 |
| Benzo(ghi)perylene | 5.829 | | ug/L | 0.48 | 10 | Į. | 58 | 30-130 | 1.7 | 20 |
| Phenanthrene | 6.259 | | ug/L | 0.07 | 10 | 6 | 33 | 30-130 | 1.6 | 20 |
| Benz(a)anthracene | 5.337 | | ug/L | 0.06 | 10 | į | 53 | 30-130 | 0.0 | 20 |
| Anthracene | 6.390 | | ug/L | 0.50 | 10 | 6 | 64 | 30-130 | 4.8 | 20 |
| Acenaphthylene | 4.296 | | ug/L | 0.30 | 10 | 4 | 13 | 30-130 | 0.0 | 20 |
| Acenaphthene | 5.210 | | ug/L | 0.50 | 10 | į | 52 | 30-130 | 0.0 | 20 |
| 2-Methylnaphthalene | 2.949 | 1 | ug/L | 0.50 | 10 | 2 | 29 | 30-130 | 18.8 | 20 |
| Fluorene | 5.519 | | ug/L | 0.50 | 10 | Ę | 55 | 30-130 | 3.6 | 20 |
| Naphthalene | 1.853 | l, r | ug/L | 0.50 | 10 | | 19 | 30-130 | 38.3 | 20 |
| Benzo(a)pyrene | 5.789 | | ug/L | 0.20 | 10 | | 58 | 30-130 | 7.1 | 20 |
| Surrogate: % 2-Fluorobiphenyl | 1.939 | | ug/L | | 5 | ; | 39 | 30-130 | | |
| Surrogate: % Terphenyl-d14 | 3.902 | | ug/L | | 5 | ; | 78 | 30-130 | | |
| Surrogate: % Nitrobenzene-d5 | 0.5641 | l, r | ug/L | | 5 | | 11 | 30-130 | | |
| MS (CC17546-MS) | | | Source: CO | 17546 | Pre | epared: 22-Dec- | 18 An | alyzed: 26-D | ec-18 | |
| Benzo(ghi)perylene | 3.234 | r | ug/L | 0.48 | 10 | | 32 | 30-130 | | 20 |
| Benzo(k)fluoranthene | 3.706 | | ug/L | 0.30 | 10 | ; | 37 | 30-130 | | 20 |
| Chrysene | 4.131 | | ug/L | 0.50 | 10 | 4 | 1 1 | 30-130 | | 20 |
| Fluoranthene | 5.232 | r | ug/L | 0.50 | 10 | į | 52 | 30-130 | | 20 |
| Indeno(1,2,3-cd)pyrene | 3.515 | r | ug/L | 0.10 | 10 | ; | 35 | 30-130 | | 20 |
| Benzo(b)fluoranthene | 3.588 | | ug/L | 0.08 | 10 | ; | 36 | 30-130 | | 20 |
| Pyrene | 5.444 | r | ug/L | 0.50 | 10 | | 54 | 30-130 | | 20 |
| Dibenz(a,h)anthracene | 4.071 | r | ug/L | 0.10 | 10 | 4 | 1 1 | 30-130 | | 20 |
| Phenanthrene | 5.461 | | ug/L | 0.07 | 10 | | 55 | 30-130 | | 20 |
| Benz(a)anthracene | 3.871 | | ug/L | 0.06 | 10 | ; | 39 | 30-130 | | 20 |
| Anthracene | 5.366 | | ug/L | 0.50 | 10 | | 54 | 30-130 | | 20 |
| Acenaphthylene | 4.862 | | ug/L | 0.30 | 10 | 4 | 19 | 30-130 | | 20 |
| Acenaphthene | 5.975 | r | ug/L | 0.50 | 10 | 6 | 30 | 30-130 | | 20 |
| 2-Methylnaphthalene | 4.284 | | ug/L | 0.50 | 10 | 4 | 13 | 30-130 | | 20 |
| Fluorene | 5.447 | r | ug/L | 0.50 | 10 | Į. | 54 | 30-130 | | 20 |
| Naphthalene | 4.516 | r | ug/L | 0.50 | 10 | 4 | 1 5 | 30-130 | | 20 |
| Benzo(a)pyrene | 3.418 | | ug/L | 0.20 | 10 | ; | 34 | 30-130 | | 20 |
| Surrogate: % 2-Fluorobiphenyl | 2.207 | r | ug/L | | 5 | 4 | 14 | 30-130 | | |
| Surrogate: % Terphenyl-d14 | 2.186 | | ug/L | | 5 | 4 | 14 | 30-130 | | |
| Surrogate: % Nitrobenzene-d5 | 1.770 | r | ug/L | | 5 | ; | 35 | 30-130 | | |
| MSD (CC17546-MSD) | | | Source: CO | C17546 | Pre | epared: 22-Dec- | 18 An | alyzed: 26-D | ec-18 | |
| Benzo(k)fluoranthene | 2.550 | | ug/L | 0.30 | 7.69231 | - | 33 | 30-130 | 11.4 | 20 |
| Chrysene | 3.328 | | ug/L | 0.50 | 7.69231 | | 13 | 30-130 | 4.8 | 20 |

| analyte(s) | Result | Flag | Units | *RDL | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit |
|-------------------------------|--------|------|------------|--------|----------------|------------------|-----------|----------------|-------|--------------|
| W8270D (SIM) | | | | | | | | | | |
| Satch 461099A - SW3520C | | | | | | | | | | |
| MSD (CC17546-MSD) | | | Source: CO | C17546 | Pre | epared: 22- | Dec-18 An | alyzed: 26-D | ec-18 | |
| Dibenz(a,h)anthracene | 2.179 | m, r | ug/L | 0.10 | 7.69231 | | 28 | 30-130 | 37.7 | 20 |
| Fluoranthene | 5.046 | r | ug/L | 0.50 | 7.69231 | | 66 | 30-130 | 23.7 | 20 |
| Fluorene | 5.185 | r | ug/L | 0.50 | 7.69231 | | 67 | 30-130 | 21.5 | 20 |
| Benzo(ghi)perylene | 1.913 | m, r | ug/L | 0.48 | 7.69231 | | 25 | 30-130 | 24.6 | 20 |
| Phenanthrene | 5.189 | | ug/L | 0.07 | 7.69231 | | 67 | 30-130 | 19.7 | 20 |
| 2-Methylnaphthalene | 3.987 | | ug/L | 0.50 | 7.69231 | | 52 | 30-130 | 18.9 | 20 |
| Indeno(1,2,3-cd)pyrene | 2.069 | m, r | ug/L | 0.10 | 7.69231 | | 27 | 30-130 | 25.8 | 20 |
| Benzo(b)fluoranthene | 2.639 | | ug/L | 0.08 | 7.69231 | | 34 | 30-130 | 5.7 | 20 |
| Benz(a)anthracene | 3.378 | | ug/L | 0.06 | 7.69231 | | 44 | 30-130 | 12.0 | 20 |
| Anthracene | 5.053 | | ug/L | 0.50 | 7.69231 | | 66 | 30-130 | 20.0 | 20 |
| Acenaphthene | 5.786 | r | ug/L | 0.50 | 7.69231 | | 75 | 30-130 | 22.2 | 20 |
| Pyrene | 5.183 | r | ug/L | 0.50 | 7.69231 | | 67 | 30-130 | 21.5 | 20 |
| Acenaphthylene | 4.635 | | ug/L | 0.30 | 7.69231 | | 60 | 30-130 | 20.2 | 20 |
| Benzo(a)pyrene | 2.425 | | ug/L | 0.20 | 7.69231 | | 32 | 30-130 | 6.1 | 20 |
| Naphthalene | 4.286 | r | ug/L | 0.50 | 7.69231 | | 56 | 30-130 | 21.8 | 20 |
| Surrogate: % Terphenyl-d14 | 2.063 | | ug/L | | 3.84615 | | 54 | 30-130 | | |
| Surrogate: % Nitrobenzene-d5 | 1.734 | r | ug/L | | 3.84615 | | 45 | 30-130 | | |
| Surrogate: % 2-Fluorobiphenyl | 2.114 | r | ug/L | | 3.84615 | | 55 | 30-130 | | |

Notes and Definitions

c1 A LCS and LCS Duplicate were performed instead of a matrix spike and matrix spike duplicate.

c3 A Blank spike was performed instead of a matrix spike

c4 A blank MS/MSD was analyzed with this batch.

1 This parameter is outside laboratory lcs/lcsd specified recovery limits.

m This parameter is outside laboratory ms/msd specified recovery limits.

r This parameter is outside laboratory rpd specified recovery limits.

dry Sample results reported on a dry weight basis

NR Not Reported

RPD Relative Percent Difference

OG The required Matrix Spike and Matrix Spike Duplicate (MS/MSD) for Oil & Grease method 1664B can only be analyzed when the client has submitted sufficient sample volume. An extra liter per MS/MSD is required to fulfill the method QC criteria. Please refer to Chain of Custody and QC Summary (MS/MSD) of the Laboratory Report to verify ample sample volume was submitted to fulfill the requirement.

pH The method for pH does not stipulate a specific holding time other than to state that the samples should be analyzed as soon as possible. For aqueous samples the 40 CFR 136 specifies a holding time of 15 minutes from sampling to analysis. Therefore all aqueous pH samples not analyzed in the field are considered out of hold time at the time of sample receipt. All soil samples are analyzed as soon as possible after sample receipt.

<u>Laboratory Control Sample (LCS)</u>: A known matrix spiked with compound(s) representative of the target analytes, which is used to document laboratory performance.

Matrix Duplicate: An intra-laboratory split sample which is used to document the precision of a method in a given sample matrix.

<u>Matrix Spike</u>: An aliquot of a sample spiked with a known concentration of target analyte(s). The spiking occurs prior to sample preparation and analysis. A matrix spike is used to document the bias of a method in a given sample matrix.

<u>Method Blank</u>: An analyte-free matrix to which all reagents are added in the same volumes or proportions as used in sample processing. The method blank should be carried through the complete sample preparation and analytical procedure. The method blank is used to document contamination resulting from the analytical process.

Method Detection Limit (MDL): The minimum concentration of a substance that can be measured and reported with 99% confidence that the analyte concentration is greater than zero and is determined from analysis of a sample in a given matrix type containing the analyte.

Reportable Detection Limit (RDL): The lowest concentration that can be reliably achieved within specified limits of precision and accuracy during routine laboratory operating conditions. For many analytes the RDL analyte concentration is selected as the lowest non-zero standard in the calibration curve. While the RDL is approximately 5 to 10 times the MDL, the RDL for each sample takes into account the sample volume/weight, extract/digestate volume, cleanup procedures and, if applicable, dry weight correction. Sample RDLs are highly matrix-dependent.

<u>Surrogate</u>: An organic compound which is similar to the target analyte(s) in chemical composition and behavior in the analytical process, but which is not normally found in environmental samples. These compounds are spiked into all blanks, standards, and samples prior to analysis. Percent recoveries are calculated for each surrogate.

<u>Continuing Calibration Verification:</u> The calibration relationship established during the initial calibration must be verified at periodic intervals. Concentrations, intervals, and criteria are method specific.



CHAIN OF CUSTODY RECORD

□ Rush TAT - Date Needed:

☑ Standard TAT - 7 to 10 business days

Special Handling:

| ulf Oil LP | ndrew Adams | ECTRUM ANALYTICAL, INC. Realiting HANIBAL TECHNOLOGY |
|-----------------------|------------------------------|---|
| Gulf Oil LP | Invoice To: Christopher Gill | Page1 of1 |
| Site Name: | Project No: | |
| Gulf Chelsea Terminal | Gulf Chelsea | All TATs subject to laboratory approval Min. 24-hr notification needed for rushes Samples disposed after 60 days unless otherwise instr |

| HANIBAL TECHNOLOGY Lepont To: Andrew Adams Gulf Oil LP 281 Eastern Ave | Invoice To: Christopher Gill Gulf Oil LP 80 William St, Suite 400 | Project No: | Samples disposed after 60 days unless otherwise instructed. Gulf Chelsea Gulf Chelsea Terminal |
|---|---|-------------------------------|--|
| Gulf Oil LP | Gulf Oil LP | Site Name: | Gulf Chelsea Terminal |
| 281 Eastern Ave | 80 William St, Suite 400 | | |
| Chelsea, MA 02150 | Wellesley, MA 02481-3705 | Location: | 281 Eastern Ave, Chelsea State: MA |
| lephone #: 617.884.5980 | | Sampler(s): | |
| oject Mgr: Andrew Adams | P.O No.: Quote/RQN: | | 0 |
| =Field Filtered 1=Na ₂ S2O ₃ 2=HCl 3=H ₂ SO ₄ 4=HNO ₃ 5=NaOH 6=Ascorbic Acid | 4=HNO; 5=NaOH 6=Ascorbic Acid | List Preservative Code below: | elow: OA/OC Reporting Notes: |
| =CH3OH 8=NaHSO ₄ 9=Deionized Water 10=H ₃ PO ₄ | 11= none 12= | | |

| | | | | | , | S | | | | | | | | | | | | | |
|--------------------|----------------|---------------|--------------|-------------------|-----------------------------|-------------|-------------------------------------|------------|--------------------|--------------------------------|---|--------------------------|--|--------------|--------------|--------------------------|--------------------------|-----------------------|--|
| | | 602 | | | | 5052773.1 | Lab ID: | 0 | | 0 =0il S0 =Soil | DW=Dinking Water | | F=Field Filtered 7=CH3OH 8=Na | Project Mgr: | Telephone #: | Chelsea | 281 Eastern Ave | Gulf Oil LP | |
| | | Trip Blank | Ouffall 003 | Outfall 003 | Outfall 003 | Outfall 003 | Sample ID: | G= Grab | X2= | SL=Sludge A=Indoor/Ambient Air | GW=Groundwater | | F=Field Filtered 1=Na ₂ S2O; 2=HCl 3=H ₂ SO ₁ 7=CH3OH 8=NaHSO ₁ 9=Defonized Water 10=H ₂ PO ₂ | Andrew Adams | 617.884.5980 | Chelsea, MA 02150 | stern Ave | LP . | Control of the Contro |
| | | | 12-19-18 131 | 12-19-18 1319 | 8-61-21 | 12-19-18 | Date: | C=Compsite | X3= | ent Air SG=Soil Gas | SW=Surface Water WW | | 4=HNO; | | | ** | | | |
| | | | 1315 | 1315 | 1315 | 1318 | Time: | | | Gas | WW=Waste Water | | 5=NaOH 6=A | P.O.No.: | ž | < | 18 | lo. | |
| | | 7 | G | G | G | G | T | ype | | | | × | 6=Ascorbic Acid | | | Wellesley, MA 02481-3705 | 80 William St, Suite 400 | Gulf Oil LP | |
| | | 75 | WS | WS | SW | WS | | atrix | | | S-100 | | Acid | | | y, MA | m St, S | P | |
| | | 2 | | ω | | | | | Vials | | | | | | |)2481-3 | uite 40 | | |
| | | | 2 | | | | | | er Glas | | Con | | | Quote/RQN: | | 3705 | ŏ | | |
| | | 1 | | | | | - | | Glass | | Containers | 1 | | QN: | | | | | |
| | | | | | | 22 | # of | Plasti | ic | | | | | | | | | | |
| | | | | _ | | | Tee | | | | | | 4.0 | | | | | | |
| | | | | | × | × | 0& | , pH | | | | 11 3 | | | | | | | |
| | | | | × | 1 | | - | | notes | for list | | 2 | List P | | - | | | | - |
| | - | | × | 1, | | | PAF | l (ben | z(a) py | re ne é | 2. | 11 | List Preservative Code below: | | Sam | Loca | | Site | |
| | | | | \vdash | | | nap | hthale | ne) | | Analysis | | ıtive C | | Sampler(s): | Location: | | Site Name | |
| | | | - | | | | | | | | | | ode bel | | - | | | | |
| _ | | | | | - | | | | ********* | | | | low: | | | 28 | | | |
| | | | | | | A | | | | | | | | | | 281 Eastern Ave, Chelsea | | | |
| | | | | | | | T | | chlori | nated | | I | | | | m Ave, | 1 | Gulf Ch | |
| | | | | | VOC | | State-s | 2 0 | | N | MA DEI | - A | * o | l e | | Chelse | | nelsea 1 | |
| napht | ben | Rec | | MTBE | s: ben | 1 | e-specifi | ☐ Tier II* | ASP A* NJ Reduced* | Standard DQA* | MA DEP MCP CAM Report? CT DPH RCP Report? | MINOHAL | VQCF | | | B | | Gulf Chelsea Terminal | |
| halene | benzene 2 µg/L | Required MLs: | | MTBE, and ethanol | zene, n | | c reportii | | g | 2 - 2 - | ιΜ Repor | Charges | Reporti | | | S | | _ | |
| naphthalene 5 µg/L | µg/L | MLs: | | thanol | VOCs: benzene, naphthalene, | | State-specific reporting standards: | Tier IV* | ASP B* | No QC | χε' □ Ω χε' □ | administratives applying | QA/QC Reporting Notes: | | | State: | | | |
| | | | * | | lene, | | rds: | | | | | - Year | | | | MA | | | |
| | <u></u> | 1 | 6 | | 0 | | | | | | 0 0 | | | | | | | | - |

3/106/4.

IR ID#

☐ Ambient ☐ Iced

↑ Refrigerated □ DI VOA Frozen

☐ Soil Jar Frozen

Condition upon receipt: Custody Seals: Present

☐ Intact ☐ Broken

Date:

LTime:

Temp °C

EDD format: E-mail to:

5

aadams@gulfoil.com, cgill@gulfoil.com, and

benzo(a)pyrene 0.1 µg/L

jennifer.atkins@aecom.com

Refinquished by;

Batch Summary

<u>1816206</u>

General Chemistry Parameters

1816206-DUP1

1816206-SRM1

1816206-SRM2

SC52773-01 (Outfall 003)

461099A

Subcontracted Analyses

CC17546-BLK

CC17546-LCS

CC17546-LCSD

CC17546-MS

CC17546-MSD

SC52773-01 (Outfall 003)

461248A

Subcontracted Analyses

CC18506-BLK

CC18506-DUP

CC18506-LCS

SC52773-01 (Outfall 003)

461279A

Subcontracted Analyses

CC19490-BLK

CC19490-LCS

CC19490-LCSD

SC52773-01 (Outfall 003)

SC52773-02 (Trip Blank)

461365A

Subcontracted Analyses

CC19445-BLK

CC19445-LCS

CC19445-LCSD

SC52773-01 (Outfall 003)

462669A

Subcontracted Analyses

CC19490-BLK

CC19490-LCS

CC19490-LCSD

CC19490-MS

CC19490-MSD

SC52773-01 (Outfall 003)

SC52773-02 (Trip Blank)